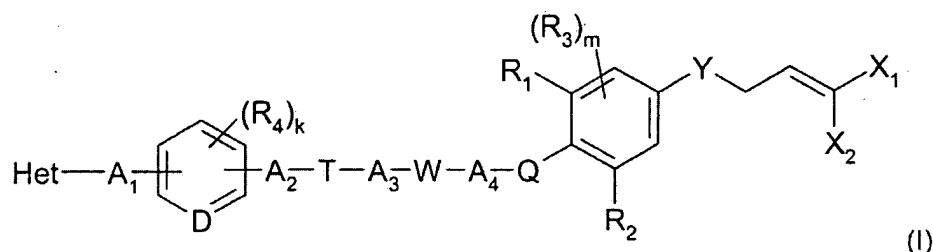


Amendments to the Claims

This listing replaces all prior versions and listing of claims in the application. Amendments are shown by addition and ~~[[deletion]]~~ or ~~deletion~~

In the Claims:

1. (Currently Amended) A compound of formula



wherein Het is a 1,2,3,4-tetrazol-5-one, which is unsubstituted or substituted with R_{iii}
~~non-aromatic heterocyclyl that does not contain cumulative double bonds and that has 5 or 6~~
~~ring members of which the linking ring member, by way of which Het is linked, by means of a~~
~~first single bond, to the remainder of the compound of formula 1, is either a nitrogen atom that~~
~~carries two further single bonds which lead to the two ring members of Het directly adjacent to~~
~~that nitrogen atom, or a carbon atom that carries a further single bond and a double bond which~~
~~lead to the two ring members of Het directly adjacent to that carbon atom, and the remaining 4~~
~~or 5 ring members of Het are, independently of one another, selected from the group consisting~~
~~of the ring members $C(R_i)(R_{ii})$, $C(=O)$, $C(=S)$, O , S , $N(R_{iii})$, $C(R_{iv})=$ and $N=$, wherein~~
~~(A) of the 5 or 6 ring members of Het, from 1 up to and including 4 ring members, independently~~
~~of one another, each contributes a hetero atom to the basic ring structure of Het consisting of 5~~
~~or 6 ring atoms, (B) two directly adjacent ring members of Het are not both O , and (C), when~~
~~the mentioned linking ring member of Het is a nitrogen atom, either (i) at least one ring member~~
~~of the mentioned remaining 4 or 5 ring members of Het is $N=$ or (ii) at least one of the 2 or 3 ring~~
~~members of Het that are neither the mentioned linking ring member of Het nor its two directly~~
~~adjacent ring members is $C(=O)$ or $C(=S)$ or (iii) at least three ring members of the~~
~~mentioned remaining 4 or 5 ring members of Het are each independently of the others $C(R_{iv})=$~~
~~or (iv) at least two ring members of the mentioned remaining 4 or 5 ring members of Het are~~
~~each independently of the other (s) O , S or $N(R_{iii})$ and, when the mentioned linking ring~~
~~member of Het is a carbon atom, either (v) the mentioned double bond starting from that carbon~~

atom leads to a nitrogen atom or (vi) the ring member of Het bonded to the mentioned further single bond starting from that carbon atom is $C(=O)$ or $C(=S)$;

R_i and R_{ii} are each independently of the other hydrogen, halogen, C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl or C_1 - C_6 alkoxy- C_1 - C_6 alkyl;

R_{iii} is C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl or C_1 - C_6 alkoxy- C_1 - C_6 alkyl;

R_{iv} is hydrogen, halogen, C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl or C_1 - C_6 alkoxy- C_1 - C_6 alkyl;

A_1 , A_2 and A_3 are each independently of the others a bond or a C_1 - C_6 alkylene bridge which is unsubstituted or substituted from one to six times by, each independently of the other(s), C_3 - C_8 cycloalkyl, C_3 - C_8 cycloalkyl- C_1 - C_6 alkyl₁ or halo- C_1 - C_3 alkyl;

A_4 is a C_1 - C_6 alkylene bridge which is unsubstituted or substituted from one to six times by, each independently of the other (s), C_3 - C_8 cycloalkyl, C_3 - C_8 cycloalkyl- C_1 - C_6 alkyl₁ or halo- C_1 - C_3 alkyl;

D is CH or N;

W is O, NR_5 , S, $S(=O)$, $S(=O)_2$, $-C(=O)-O-$, $-O-C(=O)-$, $-C(=O)-NR_6-$ or $-NR_6-C(=O)-$;

T is a bond, O, NH, NR_5 , S, $S(=O)$, $S(=O)_2$, $-C(=O)-O-$, $-O-C(=O)-$, $-C(=O)-NR_6-$ or $-NR_6-C(=O)-$;

Q is O, NR_5 , S, $S(=O)$ ₁ or $S(=O)$ ₂;

Y is O, NR_5 , S, $S(=O)$ ₁ or $S(=O)$ ₂;

X_1 and X_2 are each independently of the other fluorine, chlorine₁ or bromine;

R_1 and R_2 are each independently of the other H, halogen, CN, nitro, C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkylcarbonyl, C_2 - C_6 alkenyl, halo- C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, halo- C_2 - C_6 alkenyloxy, C_3 - C_6 alkynyloxy, C_1 - C_6 alkoxycarbonyl₁ or halo- C_3 - C_6 alkynyloxy;

R_3 is halogen, CN, nitro, C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkylcarbonyl, C_2 - C_6 alkenyl, halo- C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, halo- C_2 - C_6 alkenyloxy, C_3 - C_6 alkynyloxy, C_1 - C_6 alkoxycarbonyl₁ or halo- C_3 - C_6 alkynyloxy,

the two R_3 substituents being identical or different when m is 2;

R_4 is halogen, CN, nitro, C_1 - C_6 alkyl, halo- C_1 - C_6 alkyl, C_1 - C_6 alkylcarbonyl, C_2 - C_6 alkenyl, halo- C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, halo- C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, halo- C_2 - C_6 alkenyloxy, C_3 - C_6 alkynyloxy, C_1 - C_6 alkoxycarbonyl₁ or halo- C_3 - C_6 alkynyloxy,

the R_4 substituents being identical or different when k is greater than 1;

R₅ is H, C₁-C₆alkyl, halo-C₁-C₃alkyl, halo-C₁-C₃alkylcarbonyl, C₁-C₆alkoxyalkyl, C₁-C₆alkylcarbonyl, or C₃-C₈cycloalkyl;

R₆ is H, C₁-C₆alkyl, halo-C₁-C₃alkyl, halo-C₁-C₃alkylcarbonyl, C₁-C₆alkoxyalkyl, C₁-C₆alkylcarbonyl, or C₃-C₈cycloalkyl;

k is 0, 1, 2, or 3, when D is N_i; or

k is 0, 1, 2, 3, or 4, when D is CH; and

m is 0, 1, or 2,

or and, where applicable, possible E/Z isomers, mixtures of E/Z isomers, and/or tautomers thereof,

in each case in free form or in salt form.

2. (Original) A compound according to claim 1 in free form.
3. (Previously Presented) A compound according to claim 1, wherein X₁ and X₂ are chlorine or bromine.
4. (Previously Presented) A compound according to claim 1, wherein A₁ is a bond.
5. (Previously Presented) A compound according to claim 1, wherein the group A₂-T-A₃ is a bond.
6. (Currently Amended) A compound according to claim 1, wherein W is -O-, -C(=O)O-, or -C(=O)NH-.
7. (Previously Presented) A compound according to claim 1, wherein A₄ is a straight-chain alkylene bridge.
8. (Previously Presented) A compound according to claim 1, wherein Q is oxygen.
9. (Previously Presented) A compound according to claim 1, wherein Y is oxygen.
10. (Previously Presented) A compound according to claim 1, wherein R₁ and R₂ are bromine or chlorine.
11. (Previously Presented) A compound according to claim 1, wherein m is 0.

12. (Previously Presented) A compound according to claim 1, wherein R_4 is halogen and k is 2 or 0.

13. (Previously Presented) A compound according to claim 1, wherein D is CH.

14. (Previously Presented) A pesticidal composition comprising as active ingredient at least one compound according to claim 1, in free form or in agrochemically usable salt form, and at least one adjuvant.

15. (Original) A process for the preparation of a composition as described in claim 14, which comprises intimately mixing the active ingredient with the adjuvants.

16. (Currently Amended) A method of controlling one or more pests selected from the group consisting of insects and representatives of the order Acarina, which comprises applying a composition as described in claim 14 to the pests or to the locus thereof.